

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072221\
 Data File : BM031052.D
 Acq On : 21 Jul 2021 21:00
 Operator : CG/JU
 Sample : SST0.810
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8011

Quant Time: Jul 22 01:41:58 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 22 01:40:36 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.639	152	891	0.400	ng/ul	0.00
4) Naphthalene-d8	10.406	136	3359	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.265	164	2237	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.013	188	4961	0.400	ng/ul	0.00
17) Chrysene-d12	21.198	240	4818	0.400	ng/ul	0.00
23) Perylene-d12	23.375	264	4451	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	739	0.807	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.993	152	4792	0.820	ng/ul	0.00
18) Fluoranthene-d10	19.043	212	11888	0.733	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.251	88	801	0.784	ng/ul	92
5) Naphthalene	10.451	128	8136	0.812	ng/ul	98
7) 2-Methylnaphthalene	12.070	142	5816	0.821	ng/ul	99
8) 1-Methylnaphthalene	12.290	142	5730	0.828	ng/ul	100
10) Acenaphthylene	13.986	152	7259	0.662	ng/ul	97
11) Acenaphthene	14.330	153	6448	0.815	ng/ul	97
12) Fluorene	15.319	166	7659	0.814	ng/ul	100
14) Pentachlorophenol	16.662	266	1339	0.764	ng/ul	99
15) Phenanthrene	17.055	178	12701	0.816	ng/ul	99
16) Anthracene	17.145	178	11919	0.799	ng/ul	97
19) Fluoranthene	19.073	202	15491	0.730	ng/ul	99
20) Pyrene	19.436	202	15616	0.727	ng/ul	98
21) Benzo(a)anthracene	21.183	228	15305	0.735	ng/ul	100
22) Chrysene	21.234	228	16455	0.765	ng/ul	99
24) Benzo(b)fluoranthene	22.723	252	17077	0.793	ng/ul	95
25) Benzo(k)fluoranthene	22.767	252	17364	0.802	ng/ul	94
26) Benzo(a)pyrene	23.278	252	14665	0.784	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	25.531	276	19430	0.766	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.548	278	15592	0.760	ng/ul	96
29) Benzo(g,h,i)perylene	26.190	276	16485	0.766	ng/ul	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072221\
Data File : BM031052.D
Acq On : 21 Jul 2021 21:00
Operator : CG/JU
Sample : SST0.810
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SSTD0.8011

Quant Time: Jul 22 01:41:58 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 22 01:40:36 2021
Response via : Initial Calibration

